Package ‘agrmt’

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Type Package
Title Calculate Agreement or Consensus in Ordered Rating Scales
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Description Calculate agreement or consensus in ordered rating scales. The package implements van der Eijk's (2001) <DOI: 10.1023/A:1010374114305> measure of agreement A, which can be used to describe agreement, consensus, or polarization among respondents. It also implements measures of consensus (dispersion) by Leik, Tatsle and Wierman, Blair and Lacy, Kvalseth, Berry and Mielke, and Garcia-Montalvo and Reynal-Querol. Furthermore, an implementation of Galtung’s AJUS-system is provided to classify distributions, as well as a function to identify the position of multiple modes.

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Description

This package calculates agreement in ordered rating scales. It implements van der Eijk’s (2001) measure of agreement A, which can be used to describe agreement, consensus, or polarization among respondents. It also implements other related measures.

Details

The main functions in this package are `agreement` to calculate agreement "A", and `polarization` to calculate a polarization score based on agreement A. The package also includes functions to classify distributions according to Galtung’s (1969) AJUS-system, and changes over time according to Galtung’s (1969) ISD-system. Moreover, the function `modes` can identify the position of multiple modes.

Author(s)

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agreement

References


agreement

Calculate van der Eijk’s measure of agreement A

Description

Calculate agreement in ordered rating scales. This function implements van der Eijk’s (2001) measure of agreement A, which can be used to describe agreement, consensus, or polarization among respondents.

Usage

agreement(V, old = FALSE)

Arguments

V
A frequency vector

old
Optional argument if you wish to use the old algorithm for agreement A, as outlined in van der Eijk’s article. There is normally no reason to set the old argument.

Details

This is the main function to calculate agreement. A frequency vector describes the number of observations in a given category. For example, the vector [10,20,30,15,4] describes 10 observations with position 1, 20 observations with position 2, 30 observations with position 3, 15 observations with position 4, and 4 observations with position 5. At least three categories are required to calculate agreement.

Value

The function returns the measure of agreement A. A is 1 if there is perfect unimodality (=agreement); A is 0 if there is perfect uniformity; A is -1 if there is perfect bimodality (=lack of agreement)

Author(s)

Didier Ruedin

References

Description

Calculate agreement in ordered rating scales, but simulates coding error.

Usage

agreementError(V, n=500, e=0.01, pos=FALSE)

Arguments

V  A vector with an entry for each individual
n  Number of samples in the simulation
e  Proportion of samples for which errors are simulated
pos  Vector of possible positions. If FALSE, the values occurring in V are set as the possible values

Details

This function calculates agreement A, but simulates coding error. This can be useful to estimate standard errors and central tendency if certain positions are not observed. If all positions are observed in the vector V, bootstrapping can be used to estimate standard errors. If certain positions are not observed, bootstrapping is limited. Take an extreme example: [3 0 0 0 0]. Here we have three observations at the first position, but none at the others. Bootstrapping will always lead to the same agreement score. This can be misleading if coding error can be assumed. For example, if these three observations refer to a ‘strongly agree’ answer, it is usually conceivable that these answers could refer to ‘somewhat agree’. This function lets you specify how many of the observations should be assumed to be potentially mis-coded, and calculates agreement accordingly. If an observation is assumed to be potentially mis-coded, it is randomly set to the position to the left, the position to the right, or the position itself. If the first or last observation is chosen, the simulation takes care not to suggest values that could not occur.

You can run the function a few (hundred) times to get summary statistics of the result (mean, median, standard deviation, etc.). The function compareAgreement does just this, and compares the result with the agreement score if no coding error is assumed.
Value

The function returns the measure of agreement $A$.

Author(s)

Didier Ruedin

See Also

agreement, compareAgreement, collapse

Examples

# Sample data:
V <- c(1,1,1,2,3,3,3,4,4,4,4,4)
# Calculate agreement $A$ with coding error:
agreementError(V)
# Assume that all values could have coding error:
agreementError(V, e=1)
# Run the function a few times and show the mean:
z <- replicate(1000, agreementError(V))
  mean(z)
  hist(z) # etc.
  # you could also use the compareAgreement function.

ajas

Classify distributions

Description

Classify distributions using the AJUS-system introduced by Galtung (1969).

Usage

ajas(V, tolerance=0.1, variant="modified")

Arguments

V A frequency vector
tolerance Specify how similar values have to be to be treated as different (optional). Differences smaller than or equal to the tolerance are ignored.
variant Strict AJUS following Galtung, or modified to include F and L types (default)
Details

This function implements the AJUS-system introduced by Galtung (1969). The input is a frequency vector. Distributions are classified as A if they are unimodal with a peak in the centre, as J if they are unimodal with a peak at either end, as U if they are bimodal with a peak at both ends, and as S if they are multimodal. In addition to Galtung’s classification, the function classifies distributions as F if there is no peak and all values are more or less the same (flat). Furthermore, a distinction is drawn between J and L distributions, depending on whether they increase or decrease: J types have a peak on the right, L types have the peak on the left. The skew is given as +1 for a positive skew, as 0 for no skew, and -1 for a negative skew.

The skew is identified by comparing the sum of values left and right of the midpoint respectively. For J-type of distributions, the skew is identified on the basis of the changes between values. This way long tails cannot influence the skew, and a single peak at the left and right-hand end can be differentiated in all cases.

The aim of the AJUS system is to reduce complexity. Initially the intuition was to classify distributions on an ad-hoc basis (i.e. eye-balling). Using an algorithm is certainly more reliable, and useful if one is interested in classifying (and comparing) a large number of distributions. The argument tolerance, however is not a trivial choice and can affect results. Use the helper function ajusCheck to check sensitivity to different values of the tolerance parameter.

You can choose between a strict AJUS classification and a modified AJUSFL classification (default). The AJUS classification does not include a type for distributions without peaks (F type), and NA is returned instead. The AJUS classification does not draw a distinction between unimodal distributions with a peak at the end: the skew needs to be considered to distinguish between increasing and decreasing cases. The modified variant (default) includes the F type and the L type along with the original AJUS types.

Value

The function returns a list. The type returns a string corresponding to the pattern described by Galtung (A,J,U,S) or (F,L). The skew returns a number to describe the direction of the skew. The pattern returns the simplified pattern of the distribution. It indicates whether two values were considered the same (0), or if there was an increase (1) or decrease (-1) between two consecutive values. The length of the pattern is equal to the length of the frequency vector minus one.

Author(s)

Didier Ruedin

References


See Also

isd, ajusCheck, ajusPlot
Sensitivity test for AJUS

**Description**

Check sensitivity of AJUS to different tolerance parameters.

**Usage**

```
ajusCheck(V, t=seq(from=0.05, to=0.2, by=0.05), variant="modified")
```

**Arguments**

- `V`: A frequency vector
- `t`: A vector of tolerance parameters to check. Differences smaller than or equal to the tolerance are ignored.
- `variant`: Strict AJUS following Galtung, or modified to include F and L types (default)

**Details**

This function runs the AJUS system with a range of tolerance parameters. You can easily check how sensitive the classification of the distribution is to the tolerance parameter.

**Value**

The function returns a list. The `tolerance` returns the tolerance parameters tested. The `type` returns a series of strings corresponding to the pattern described by Galtung (A,J,U,S) or (F, L) for each tolerance parameter. The `skew` returns a number to describe the direction of the skew. See `ajus` for a description of the different arguments and the AJUS types.

**Author(s)**

Didier Ruedin

**See Also**

`ajus`
ajusPlot  

Plot vector with AJUS type

Description

Plot a frequency vector among with its AJUS type.

Usage

    ajusPlot(V, tolerance=0.1, variant="modified", ...)

Arguments

V  A frequency vector

tolerance  Specify how similar values have to be to be treated as different (optional). Differences smaller than or equal to the tolerance are ignored.

...  Arguments to pass to the plotting function

variant  Strict AJUS following Galtung, or modified to include F and L types (default)

Details

This function plots the frequency vector along with its AJUS classification and skew. See ajus for a description of the AJUS system and the different parameters. In contrast to the ajus function, ajusPlot can deal with missing values (they are removed when calculating the AJUS type, but considered in the plot). This makes ajusPlot useful for classifying time series. Additional arguments can be passed to the to the underlying plot function.

Author(s)

Didier Ruedin

See Also

ajus

BerryMielke  

Calculate IOV

Description

Calculate Berry and Mielke's IOV.

Usage

    BerryMielke(V)
Arguments

V  A frequency vector

Details

This function calculates Berry and Mielke’s IOV, a measure of dispersion based on squared Euclidean distances. This function follows the presentation by Blair and Lacy 2000, but includes the adjustment for Tmax omitted by Blair and Lacy as there is no reason to leave it out. The derived measure COV by Kvalseth is implemented as Kvalseth. Usually, the IOV is equivalent to 1-lsquared.

Value

The function returns the IOV.

Author(s)

Didier Ruedin

References


See Also

lsquared, Kvalseth

Examples

# Sample data
V <- c(30, 40, 210, 130, 530, 50, 10)
BerryMielke(V)

BlairLacy  Calculate l

Description

Calculate Blair and Lacy’s l.

Usage

BlairLacy(V)
Arguments
V A frequency vector

Details
This function calculates Blair and Lacy’s l, a measure of concentration based on linear Euclidean distances. This function follows the presentation by Blair and Lacy 2000. The measure l-squared by Blair and Lacy is implemented as lsquared.

Value
The function returns the l.

Author(s)
Didier Ruedin

References

See Also
lsquared

Examples
# Sample data
V <- c(30,40,210,130,530,50,10)
BlairLacy(V)

collapse Reduces a vector to a frequency vector

Description
This function reduces a vector to a frequency vector.

Usage
collapse(D,pos=FALSE)

Arguments
D Vector
pos Optional: position of categories
Details

This function reduces a vector to a frequency vector. This function is similar to the way `table` summarizes vectors, but this function can deal with categories of frequency 0 (if the argument `pos` is specified). Here we assume a vector with an entry for each individual (sorted in any way). Each entry states the position of an individual. When the number of positions is naturally limited, such as when categorical positions are used, frequency vectors can summarize this information: how many individuals have position 1, how many individuals have position 2, etc. A frequency vector has an entry for each position in the population (sorted in ascending order). Each entry states the number of individuals in the population with this position.

The argument `pos` is required if certain positions do not occur in the population. For example, if we have positions on a 7-point scale, and position 3 never occurs in the population, the argument `pos` must be specified. In this case, our argument may be `pos=1:7`. We can also use categories more generally, as in `c(-3,-1,0,.5,1,2,5)`. Specifying the positions of categories when all positions actually occur in the population has no side-effects.

Value

A frequency vector

Author(s)

Didier Ruedin

See Also

expand

**compareAgreement**  
*Compare agreement A with and without simulated coding error*

Description

Calculate agreement in ordered rating scales, and compares this to agreement with simulated coding error.

Usage

```r
compareAgreement(V, n=500, e=0.01, N=500, pos=FALSE)
```

Arguments

- `V`  
  A vector with an entry for each individual
- `n`  
  Number of samples in the simulation of coding errors
- `e`  
  Proportion of samples for which errors are simulated
- `N`  
  Number of replications for calculating mean and standard deviation
- `pos`  
  Vector of possible positions. If FALSE, the values occurring in V are set as the possible values
Details

This function calculates agreement on a vector, and compares the value with agreement with simulated coding error. It runs the function `agreementError` N times. The other arguments (n, e, pos) are passed down to the `agreementError` function.

Value

The function returns a list with agreement A without simulated coding errors, the mean of agreement with simulated coding error, and the standard deviation of agreement with simulated coding error.

Author(s)

Didier Ruedin

See Also

`agreement`, `agreementError`

Examples

```r
# Sample data:
V <- c(1,1,1,2,3,3,3,4,4,4,4,4)
compareAgreement(V)
```

**compareValues**

*Compares two values*

Description

This is a helper function to compare two values.

Usage

`compareValues(A,B,tolerance=0.1)`

Arguments

- **A**: A number
- **B**: A number
- **tolerance**: Specify how similar values have to be to be treated as different. Differences smaller than or equal to the tolerance are ignored.

Details

This is a helper function compare two values. Two values are more or less the same, or one of the two is bigger.
Value

The function returns number to describe the relationship: -1 if A is bigger, 1 if B is bigger, and 0 if the two are more or less the same.

Author(s)

Didier Ruedin

Description

Calculate consensus in ordered rating scales. This function implements Tastle and Wierman’s (2007) measure of consensus (ordinal dispersion), which can be used to describe agreement, consensus, dispersion, or polarization among respondents.

Usage

consensus(v)

Arguments

v A frequency vector

Details

This function calculates consensus following Tastle and Wierman (2007). The measure of consensus is based on the Shannon entropy. A frequency vector describes the number of observations in a given category. For example, the vector [10,20,30,15,4] describes 10 observations with position 1, 20 observations with position 2, 30 observations with position 3, 15 observations with position 4, and 4 observations with position 5.

If you come across an error that the vector supplied does not contain whole numbers, try round(v, 0) to remove any detritus from calculating the frequency vector.

Value

The function returns the measure of consensus. It is 1 if there is perfect uniformity; it is 0 if there is perfect bimodality (=lack of agreement)

Author(s)

Didier Ruedin

References

dsquared

See Also

agreement

Examples

# Sample data
V <- c(30, 40, 210, 130, 530, 50, 10)
# Calculate consensus
consensus(V)
# The degree of consensus is given as 0.7256876

dSquared |
---
Calculate d-squared

Description

Calculate Blair and Lacy’s d-squared.

Usage

dSquared(V)

Arguments

V | A frequency vector

Details

This function calculates Blair and Lacy’s d-squared, a measure of concentration based on squared Euclidean distances. This function follows the presentation by Blair and Lacy 2000. The measure l-squared normalizes the values and is implemented as lsquared.

Value

The function returns the d-squared.

Author(s)

Didier Ruedin

References


See Also

lsquared, BlairLacy
**entropy**

**Examples**

```
# Sample data
V <- c(30, 40, 210, 130, 530, 50, 10)
dsquared(V)
```

<table>
<thead>
<tr>
<th>entropy</th>
<th>Calculate Shannon entropy</th>
</tr>
</thead>
</table>

**Description**

Calculate Shannon entropy, following Tastle and Wierman.

**Usage**

`entropy(V)`

**Arguments**

- `V` A frequency vector

**Details**

This function calculates the Shannon entropy following Tastle and Wierman (2007). A frequency vector describes the number of observations in a given category. For example, the vector `[10, 20, 30, 15, 4]` describes 10 observations with position 1, 20 observations with position 2, 30 observations with position 3, 15 observations with position 4, and 4 observations with position 5.

This function follows Tastle and Wierman and ignores categories with zero observations. This does not follow the formula indicated.

See `consensus` for a function that considers the order of categories.

**Value**

The function returns the Shannon entropy.

**Author(s)**

Didier Ruedin

**References**


**See Also**

`consensus`
**Examples**

```
# Sample data
V <- c(30, 40, 210, 130, 530, 50, 10)
# Calculate entropy
entropy(V)
```

---

**Description**

This function expands a frequency vector to a vector.

**Usage**

```
expand(F)
```

**Arguments**

- **F** Frequency vector

**Details**

This function takes a frequency vector and expands it to a longer vector with one entry for each observation. It reverses the `collapse` function. A frequency vector has an entry for each position in the population. Each entry states the number of individuals in the population with this position. Here we create a vector with an entry for each individual.

**Value**

A vector

**Author(s)**

Didier Ruedin

**See Also**

`collapse`
Description

Classify changes over time using the ISD-system introduced by Galtung (1969).

Usage

isd(V, tolerance=0.1)

Arguments

V  A vector with length 3
tolerance  Specify how similar values have to be to be treated as different (optional). Differences smaller than or equal to the tolerance are ignored.

Details

This function implements the ISD-system introduced by Galtung (1969). The input is a vector of length 3. Each value stands for a different point in time. The ISD-system examines the two transition points, and classifies the changes over time.

Value

The function returns a list. The type returns a number corresponding to the pattern described by Galtung. The description returns a string where the two transitions are spelled out (increase, flat, decrease).

Author(s)

Didier Ruedin

References


See Also

ajus
**Kvalseth**

*Calculate Kvalseth’s COV*

---

**Description**

Calculate Kvalseth’s COV.

**Usage**

`Kvalseth(v)`

**Arguments**

- `v` A frequency vector

**Details**

This function calculates Kvalseth’s COV, a measure of dispersion based on linear Euclidean distances. It is based on the IOV measure, implemented as `BerryMielke`. This function follows the presentation by Blair and Lacy 2000.

**Value**

The function returns the COV.

**Author(s)**

Didier Ruedin

**References**


**See Also**

`BerryMielke`, `lsquared`

**Examples**

```r
# Sample data
V <- c(30, 40, 210, 130, 530, 50, 10)
Kvalseth(V)
```
Leik

Calculate ordinal dispersion

Description

Calculates ordinal dispersion as introduced by Leik (1966)

Usage

leik(v)

Arguments

v
A frequency vector

Details

This function calculates ordinal dispersion as introduced by Robert K. Leik (1966). It uses the cumulative frequency distribution to determine ordinal dispersion. The extremes (agreement, polarization) largely correspond to the types used by Cees van der Eijk. By contrast, the mid-point depends on the number of categories: it tends toward 0.5 as the number of categories increases. Leik defends this difference by highlighting the increased probability of falling into polarized patterns when there are fewer categories. If all observations are in the same category, ordinal dispersion is 0. With half the observations in one extreme category, and half the observations in the other extreme, Leik’s measure gives a value of 1.

The dispersion measure is a percentage, and can be interpreted accordingly. Ordinal dispersion can be used to express consensus or agreement, simply by taking: 1 - ordinal dispersion.

Value

The function returns the ordinal dispersion

Author(s)

Didier Ruedin

References


See Also

polarization, agreement
Examples

# Example 1:
V <- c(30,40,210,130,530,50,10)
# Calculate polarization
Leik(V)
# The ordinal dispersion is given as 0.287
polarization(V)
# Polarization is given as 0.194 (as contrast)

lsquared

Calculate l-squared

Description

Calculate Blair and Lacy’s l-squared.

Usage

lsquared(V)

Arguments

V A frequency vector

Details

This function calculates Blair and Lacy’s l-squared, a measure of concentration based on squared Euclidean distances. This function follows the presentation by Blair and Lacy 2000. The measure ‘l’ by Blair and Lacy is implemented as BlairLacy.

Value

The function returns the l-squared.

Author(s)

Didier Ruedin

References


See Also

BerryMielke, BlairLacy
minnz

Examples

# Sample data
V <- c(30, 40, 210, 130, 530, 50, 10)
lsquared(V)

<table>
<thead>
<tr>
<th>minnz</th>
<th>Non-zero minimum</th>
</tr>
</thead>
</table>

Description

Helper function to calculate the smallest value of the vector except for 0 (non-zero minimum).

Usage

minnz(V)

Arguments

V       A vector

Details

This is a helper function to calculate the non-zero minimum of a vector. The result is the smallest value of the vector, but cannot be zero.

Value

The function returns the non-zero minimum

Author(s)

Didier Ruedin

modes

Identify multiple modes

Description

Identifies (multiple) modes in a frequency vector.

Usage

modes(V, pos=FALSE, tolerance=0.1)
Arguments

- **v** A frequency vector
- **pos** Categories of frequency vector (optional)
- **tolerance** Specify how similar values have to be to be treated as different (optional). Differences smaller than or equal to the tolerance are ignored.

Details

This function identifies which positions of a frequency vector correspond to the mode. If there are multiple modes of the same value, all matching positions will be reported. Use the function `collapse` to create frequency vectors if necessary.

Value

The function returns a list. The `at` returns the categories of the frequency vector. Either these categories were specified using the argument `pos`, or we assume it to be 1:k (with k the number of categories in the frequency vector). If the length of the `pos` argument does not match the length of the frequency vector, a warning is shown, and the `pos` argument is ignored. The `frequencies` returns the frequency vector. The `mode` returns the value of the mode(s). If there are multiple modes, they are listed. Similar frequencies are counted as equal, using the tolerance argument. To prevent similar frequencies to be considered the same, set `tolerance` to 0. The `positions` returns the positions of the vector that correspond to the mode. This will differ from the `mode` if `pos` is provided. The `contiguous` returns TRUE if all modes are contiguous, and FALSE if there are different values in between. If there is only one mode, it is defined as contiguous (i.e. TRUE).

Author(s)

Didier Ruedin

See Also

- `secondModes`

Examples

```r
# Example 1: finding the mode
V1 <- c(30, 40, 210, 130, 530, 50, 10)
modes(V1) # will find position 5
# Example 2:
V2 <- c(3, 0, 4, 1)
modes(V2) # will find position 3
# Example 3: providing categories
modes(V2, pos=1:2) # will still find position 3, but give the value of 1 as mode
# Example 4: similar values
V3 <- c(30, 40, 500, 130, 530, 50, 10)
modes(V3, tolerance=30) # will find positions 3 and 5 (500 and 530 are nearly the same)
```
MRQ

Calculates MRQ polarization index

Description

This function calculates the MRQ polarization index from a population vector.

Usage

\texttt{mrq(z)}

Arguments

\texttt{z} \\
(Standardized) frequency vector

Details

This function implements the polarization index introduced by Garcia-Montalvo and Reynal-Querol (2005), also known as the Reynal-Querol index of polarization (RQ). It is a measure of dispersion based on squared Euclidean distances. The frequency vector needs to be standardized for the Reynal-Querol index to work; if the sum of the frequency vector is not 1 (i.e. it is not standardized), the function automatically standardizes the frequency vector by dividing each element of the vector by the sum of the vector. The assumption is that the frequencies are complete.

Value

Index of polarization (RQ).

Author(s)

Didier Ruedin

References


Examples

\begin{verbatim}
# Sample data
V <- c(30, 40, 210, 130, 530, 50, 10)
MRQ(V)
\end{verbatim}
patternAgreement  *Calculates patterns agreement*

**Description**
Helper function to calculate agreement A from a pattern vector.

**Usage**

```r
patternAgreement(P, old=FALSE)
```

**Arguments**

- `P`  
  A pattern vector

- `old`  
  Optional argument if the old algorithm for agreement A is to be used. There is normally no reason to set the `old` argument.

**Details**
This is a helper function to calculate agreement A from a pattern vector.

**Value**
The function returns the measure of agreement A

**Author(s)**
Didier Ruedin

**See Also**

- `agreement`

patternVector  *Creates pattern vector*

**Description**
Helper function to create a pattern vector from a frequency vector.

**Usage**

```r
patternVector(V)
```

**Arguments**

- `V`  
  A frequency vector
Details
This is a helper function to create a pattern vector from a frequency vector. A pattern vector reduced all values greater or equal to 1 to 1, and values of 0 remain 0. A frequency vector \((0,0,18,59,0,34,2)\) is turned into a pattern vector \((0,0,1,1,0,1,1)\).

Value
The function returns a pattern vector.

Author(s)
Didier Ruedin

See Also
agreement

Description
Calculates polarization, based on measure of agreement A

Usage
\[
polarization(V, old = FALSE)\]

Arguments
- **V** A frequency vector
- **old** Specify old=TRUE to use the depreciated algorithm for agreement A

Details
This function calculates polarization by re-scaling agreement A introduced by Cees van der Eijk. Whereas agreement A ranges from -1 to 1, polarization ranges from 0 to 1. If all observations are in the same category, polarization is 0. With half the observations in one category, and half the observations in a different (non-neighbouring) category, polarization is 1. Polarization is 0.5 for a uniform distribution over all categories.

Value
The function returns a polarization score

Author(s)
Didier Ruedin
See Also

agreement

Examples

V <- c(30,40,210,130,530,50,10)
# Calculate polarization
polarization(V)
# The rate of polarization is given as 0.1943333

---

reduceVector   Remove zeros and repeated values

Description

This is a helper function to remove all zeros and repeated values from a vector.

Usage

reduceVector(X)

Arguments

X   A (frequency) vector

Details

This is a helper function to strip all zeros and repeated values from a vector.

Value

The function returns vector

Author(s)

Didier Ruedin

See Also

agreement
secondModes

Most common and second most common values

Description
Identifies the most common (multiple) modes for frequency vectors as well as the second most common values.

Usage
secondModes(V, pos=FALSE, tolerance=0.1)

Arguments
- V: A frequency vector
- pos: Categories of frequency vector (optional)
- tolerance: Specify how similar values have to be to be treated as different (optional). Differences smaller than or equal to the tolerance are ignored.

Details
This function identifies which positions of a frequency vector correspond to the mode(s) as implemented in the modes function. It also reports the second most common position in the same manner.

Value
The function returns a list for the most common and the second most common value(s). The output corresponds to that of the modes function.

Author(s)
Didier Ruedin

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